

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120620\
 Data File : BN012920.D
 Acq On : 06 Dec 2020 16:41
 Operator : CG/JU
 Sample : L4859-04
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B94

Manual Integrations
 APPROVED

mohammad
 12/7/2020 2:11:41 PM

Quant Time: Dec 07 03:26:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 07 03:06:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1678	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	7003	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.06	164	4033	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.81	188	8979	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	7416	0.40	ng/ul	0.00
20) Perylene-d12	23.14	264	6330	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	1038	0.10	ng/ul	-0.02
14) Fluoranthene-d10	18.86	212	2458	0.10	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	1426	0.070	ng/ul	98
5) 2-Methylnaphthalene	11.84	142	783	0.059	ng/ul	96
7) Acenaphthylene	13.77	152	5077	0.266	ng/ul	96
8) Acenaphthene	14.12	153	2847	0.182	ng/ul	100
9) Fluorene	15.12	166	3601	0.200	ng/ul	99
12) Phenanthrene	16.85	178	102153	3.393	ng/ul	96
13) Anthracene	16.95	178	22058	0.861	ng/ul	95
15) Fluoranthene	18.89	202	323525	9.043	ng/ul	96
17) Pyrene	19.25	202	271818	7.760	ng/ul	98
18) Benzo(a)anthracene	21.02	228	158618	5.170	ng/ul	99
19) Chrysene	21.07	228	162760	4.801	ng/ul	98
21) Benzo(b)fluoranthene	22.52	252	207427m	7.083	ng/ul	
22) Benzo(k)fluoranthene	22.56	252	65211m	2.015	ng/ul	
23) Benzo(a)pyrene	23.05	252	128324m	4.723	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.20	276	86493	2.423	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.21	278	21419	0.759	ng/ul	95
26) Benzo(a,h,i)perylene	25.83	276	72779	2.254	ng/ul	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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